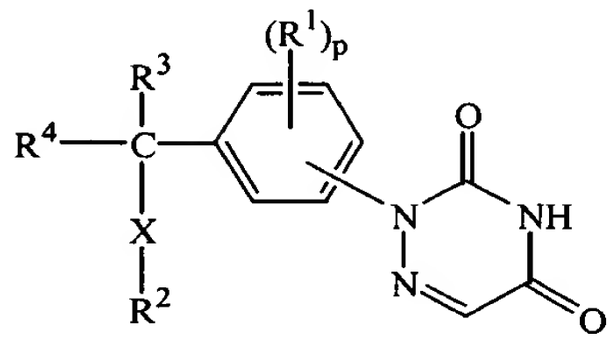
Listing of Claims

1-110. (canceled)

111. (new) A compound having the formula



(I)

a *N*-oxide, a pharmaceutically acceptable addition salt or a stereochemically isomeric form thereof, wherein:

p represents an integer being 0, 1, 2, 3 or 4;

X represents O, S, NR^5 or a direct bond;

Y represents O, S, NR^5, or $S(O)_2$;

each R^1 independently represents C_{1-6}alkyl, halo, polyhaloC_{1-6}alkyl, hydroxy, mercapto, C_{1-6}alkyloxy, C_{1-6}alkylthio, C_{1-6}alkylcarbonyloxy, aryl, cyano, nitro, Het^3, R^6, NR^7R^8 or C_{1-4}alkyl substituted with Het^3, R^6 or NR^7R^8;

R^2 represents Het^1 or C_{1-6}alkyl substituted with one or two substituents selected from hydroxy, cyano, amino, mono- or di(C_{1-4}alkyl)amino, C_{1-6}alkyloxy, C_{1-6}alkylsulfonyloxy, C_{1-6}alkyloxycarbonyl, C_{3-7}cycloalkyl, aryl, aryloxy, arylthio, Het^1, Het^1oxy, and Het^1thio; and if X is O, S or NR^5, then R^2 may also represent aminocarbonyl, aminothiocarbonyl, C_{1-4}alkylcarbonyl, C_{1-4}alkylthiocarbonyl, arylcarbonyl, arylthiocarbonyl, Het^1carbonyl or Het^1thiocarbonyl;

R^3 represents hydrogen, C_{1-6}alkyl or C_{3-7}cycloalkyl;

R^4 represents hydrogen, C_{1-6}alkyl or C_{3-7}cycloalkyl; or

R^3 and R^4 taken together form a C_{2-6}alkanediyl;

R^5 represents hydrogen or C_{1-4}alkyl;

each R^6 independently represents C_{1-6}alkylsulfonyl, aminosulfonyl, mono- or di(C_{1-4}alkyl)aminosulfonyl, mono- or di(benzyl)aminosulfonyl, polyhaloC_{1-6}alkylsulfonyl, C_{1-6}alkylsulfinyl, phenylC_{1-4}alkylsulfonyl, piperazinylsulfonyl, aminopiperidinylsulfonyl, piperidinylaminosulfonyl, *N*-C_{1-4}alkyl-*N*-piperidinylaminosulfonyl or mono-or di(C_{1-4}alkyl)aminoC_{1-4}alkylsulfonyl;

each R^7 and each R^8 are independently selected from the group consisting of: hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, aryl, arylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, aminocarbonyl, arylcarbonyl, Het^3carbonyl, C_{1-4}alkylcarbonyloxy-C_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)aminoC_{1-4}alkyl, arylaminocarbonyl, arylaminothiocarbonyl, Het^3aminocarbonyl, Het^3aminothiocarbonyl, C_{3-7}cycloalkyl, pyridinylC_{1-4}alkyl, C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3, Het^4 and R^6;

R^9 and R^{10} are each independently selected from the group consisting of: hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, phenyl, phenylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, aminocarbonyl, phenylcarbonyl, Het^3carbonyl, C_{1-4}alkylcarbonyloxy-C_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)aminoC_{1-4}alkyl, phenylaminocarbonyl, phenylaminothiocarbonyl, Het^3aminocarbonyl, Het^3aminothiocarbonyl, C_{3-7}cycloalkyl, pyridinylC_{1-4}alkyl, C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3, Het^4 and R^6;

each R^{11} independently being selected from the group consisting of: hydroxy, mercapto, cyano, nitro, halo, trihalomethyl, C_{1-4}alkyloxy, formyl, trihaloC_{1-4}alkylsulfonyloxy, R^6, NR^7R^8, C(=O)NR^7R^8, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, aryl, aryloxy, arylcarbonyl, C_{3-7}cycloalkyl, C_{3-7}cycloalkyloxy, phthalimide-2-yl, Het^3 and C(=O)Het^3;

R^{12} and R^{13} are each independently selected from the group consisting of: hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, phenyl, phenylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, phenylcarbonyl, C_{1-4}alkylcarbonyloxyC_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)aminoC_{1-4}alkyl, phenylaminocarbonyl, phenylaminothiocarbonyl, C_{3-7}cycloalkyl, pyridinylC_{1-4}alkyl, C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14} and R^6;

each R^{14} independently represents hydrogen, C_{1-4}alkyl, C_{3-7}cycloalkyl, aminocarbonylmethylene or mono-or di(C_{1-4}alkyl)aminocarbonylmethylene;

aryl represents phenyl optionally substituted with one, two or three substituents each independently selected from nitro, azido, cyano, halo, hydroxy, C_{1-4}alkyl, C_{3-7}cycloalkyl, C_{1-4}alkyloxy, formyl, polyhaloC_{1-4}alkyl, NR^9R^{10}, C(=O)NR^9R^{10}, C(=O)-O-R^{14}, R^6, -O-R^6, phenyl, Het^3, C(=O)Het^3 and C_{1-4}alkyl substituted with hydroxy, C_{1-4}alkyloxy, C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3 or NR^9R^{10};

Het^1 represents a heterocycle selected from the group consisting of: pyrrolyl, pyrrolinyl, imidazolyl, imidazolinyl, pyrazolyl, pyrazolinyl, triazolyl, tetrazolyl, furanyl, tetrahydrofuranyl, thienyl, thiolanyl, dioxolanyl, oxazolyl, oxazolinyl, isoxazolyl, thiazolyl, thiazolinyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl, pyrrolidinyl, piperidinyl, piperazinyl, morpholinyl, thiomorpholinyl, dioxanyl, dithianyl, trithianyl, triazinyl, benzothienyl, isobenzothienyl, benzofuranyl, isobenzofuranyl, benzothiazolyl, benzoxazolyl, indolyl, isoindolyl, indolinyl, purinyl, 1*H*-pyrazolo[3,4-d]pyrimidinyl, benzimidazolyl, quinolyl, isoquinolyl, cinnolinyl, phtalazinyl, quinazolinyl, quinoxalinyl, thiazolopyridinyl, oxazolopyridinyl and imidazo[2,1-b]thiazolyl; wherein said heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with one or two substituents independently selected from Het^2 and R^{11}; provided Het^1 is other than 2-substituted-pyridin-5-yl;

Het^2 represents a heterocycle selected from the group consisting of: pyrrolyl, pyrrolinyl, imidazolyl, imidazolinyl, pyrazolyl, pyrazolinyl, triazolyl, tetrazolyl, furanyl, tetrahydrofuranyl, thienyl, thiolanyl, dioxolanyl, oxazolyl, oxazolinyl, isoxazolyl, thiazolyl, thiazolinyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl, dioxanyl, dithianyl, trithianyl, triazinyl, benzothienyl, isobenzothienyl, benzofuranyl, isobenzofuranyl, benzothiazolyl, benzoxazolyl, indolyl, isoindolyl, indolinyl, purinyl, 1*H*-pyrazolo[3,4-d]pyrimidinyl, benzimidazolyl, quinolyl, isoquinolyl, cinnolinyl, phtalazinyl, quinazolinyl, quinoxalinyl, thiazolopyridinyl, oxazolopyridinyl and imidazo[2,1-b]thiazolyl; wherein said heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^4, R^{11} and C_{1-4}alkyl optionally substituted with one or two substituents independently selected from Het^4 and R^{11};

Het^3 represents a monocyclic heterocycle selected from the group consisting of:\ pyrrolidinyl, piperidinyl, piperazinyl, morpholinyl, thiomorpholinyl and tetrahydropyranyl; wherein said monocyclic heterocycles each independently may optionally be substituted with, where possible, one, two, three or four substituents each independently selected from hydroxy, C_{1-4}alkyl, C_{1-4}alkyloxy, C_{1-4}alkylcarbonyl, piperidinyl, $NR^{12}R^{13}$, C(=O)-O-R^{14}, R^6 and C_{1-4}alkyl substituted with one or two substituents independently selected from hydroxy, C_{1-4}alkyloxy, phenyl, C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, R^6 and $NR^{12}R^{13}$;

Het^4 represents a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl,

isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl.

112. (new) A compound of claim 111 wherein:

each R^7 and each R^8 are independently selected from the group consisting of: hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, aryl, arylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, aminocarbonyl, arylcarbonyl, Het^3carbonyl, C_{1-4}alkylcarbonyloxy-C_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)aminoC_{1-4}alkyl, arylaminocarbonyl, arylaminothiocarbonyl, Het^3aminocarbonyl, Het^3aminothiocarbonyl, C_{3-7}cycloalkyl, pyridinylC_{1-4}alkyl, C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3 and R^6;

R^9 and R^{10} are each independently selected from the group consisting of: hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, phenyl, phenylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, aminocarbonyl, phenylcarbonyl, Het^3carbonyl, C_{1-4}alkylcarbonyloxy-C_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)aminoC_{1-4}alkyl, phenylaminocarbonyl, phenylaminothiocarbonyl, Het^3aminocarbonyl, Het^3aminothiocarbonyl, C_{3-7}cycloalkyl, pyridinylC_{1-4}alkyl, C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3 and R^6;

R^{11} is being selected from the group consisting of: hydroxy, mercapto, cyano, nitro, halo, trihalomethyl, C_{1-4}alkyloxy, formyl, trihaloC_{1-4}alkylsulfonyloxy, R^6, NR^7R^8, $C(=O)NR^7R^8$, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, aryl, aryloxy, arylcarbonyl, C_{3-7}cycloalkyl, C_{3-7}cycloalkyloxy, phthalimide-2-yl, Het^3, Het^4 and C(=O)Het^3; and

Het^2 represents a heterocycle selected from the group consisting of: pyrrolyl, pyrrolinyl, imidazolyl, imidazolinyl, pyrazolyl, pyrazolinyl, triazolyl, tetrazolyl, furanyl, tetrahydrofuranyl, thienyl, thiolanyl, dioxolanyl, oxazolyl, oxazolinyl, isoxazolyl, thiazolyl, thiazolinyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl, dioxanyl, dithianyl, trithianyl, triazinyl, benzothienyl, isobenzothienyl, benzofuranyl, isobenzofuranyl, benzothiazolyl, benzoxazolyl, indolyl, isoindolyl, indolinyl, purinyl, 1*H*-pyrazolo[3,4-d]pyrimidinyl, benzimidazolyl, quinolyl, isoquinolyl, cinnolinyl, phtalazinyl, quinazolinyl, quinoxalinyl, thiazolopyridinyl, oxazolopyridinyl and imidazo[2,1-b]thiazolyl; wherein said heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently

selected from R^{11} and C_{1-4}alkyl optionally substituted with one or two substituents independently selected from R^{11}.

113. (new) A compound of the formula:
2-[3,5-dichloro-4-[1-methyl-1-(4-phenyl-2-thiazolyl)ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;
2-[3,5-dichloro-4-[1-[4-(3-chlorophenyl)-5-methyl-2-thiazolyl]-1-methylethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;
2-[3,5-dichloro-4-[1-methyl-1-(5-phenyl-1,2,4-oxadiazol-3-yl)ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;
2-[3,5-dichloro-4-[1-(4,5-diphenyl-2-thiazolyl)-1-methylethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;
2-[3,5-dichloro-4-[1-methyl-1-[5-(2-methylphenyl)-1,2,4-oxadiazol-3-yl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;
2-[3,5-dichloro-4-[1-methyl-1-(4-methyl-5-phenyl-2-thiazolyl)ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;
2-[3,5-dichloro-4-[1-methyl-1-[4-phenyl-5-(3-pyridinyl)-2-thiazolyl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;
2-[3,5-dichloro-4-[1-methyl-1-[4-phenyl-5-(phenylmethyl)-2-thiazolyl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;
2-[3,5-dichloro-4-[1-methyl-1-[5-(4-pyridinyl)-1,2,4-oxadiazol-3-yl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;
2-[3,5-dichloro-4-[1-methyl-1-[4-(3-thienyl)-2-thiazolyl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;
2-[3,5-dichloro-4-[1-[4-(2-furanyl)-2-thiazolyl]-1-methylethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;
2-[3,5-dichloro-4-[1-methyl-1-[5-(3-pyridinyl)-1,2,4-oxadiazol-3-yl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione;
2-[3,5-dichloro-4-[1-methyl-1-[5-(2-methyl-3-pyridinyl)-1,2, 4-oxadiazol-3-yl]ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione; or
2-[3,5-dichloro-4-[1-methyl-1-(5-phenyl-1,3,4-oxadiazol-2-yl)ethyl]phenyl]-1,2,4-triazine-3,5(2*H*,4*H*)-dione; or a *N*-oxide, a pharmaceutically acceptable addition salt or a stereochemically isomeric form thereof.

114. (new) A composition comprising a pharmaceutically acceptable carrier and, as active ingredient, a therapeutically effective amount of a compound of claim 111.

115. (new) A compound of claim 111 provided that in those compounds wherein X is a direct bond, at least one of R^3 and R^4 is hydrogen, and R^2 is 3-pyridinyl optionally substituted in the 6 position with an optionally substituted alkyl or acyl group are excluded.

116. (new) A compound of claim 112 provided that in those compounds wherein X is a direct bond, at least one of R^3 and R^4 is hydrogen, and R^2 is 3-pyridinyl optionally substituted in the 6 position with an optionally substituted alkyl or acyl group are excluded.

117. (new) A compound of claim 111 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

118. (new) A compound of claim 112 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

119. (new) A compound of claim 115 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

120. (new) A compound of claim 116 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

121. (new) A compound of claim 111 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

122. (new) A compound of claim 112 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

123. (new) A compound of claim 115 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl,

thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

124. (new) A compound of claim 116 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

125. (new) A compound of claim 117 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

126. (new) A compound of claim 118 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

127. (new) A compound of claim 119 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

128. (new) A compound of claim 120 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

129. (new) A compound of claim 111 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

130. (new) A compound of claim 112 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

131. (new) A compound of claim 115 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

132. (new) A compound of claim 116 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

133. (new) A compound of claim 117 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

134. (new) A compound of claim 118 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

135. (new) A compound of claim 119 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

136. (new) A compound of claim 120 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

137. (new) A compound of claim 121 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

138. (new) A compound of claim 122 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

139. (new) A compound of claim 123 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

140. (new) A compound of claim 124 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

141. (new) A compound of claim 125 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

142. (new) A compound of claim 126 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

143. (new) A compound of claim 127 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

144. (new) A compound of claim 128 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

145. (new) A compound of claim 111 wherein p is 1 or 2 and each R^1 is chloro.

146. (new) A compound of claim 112 wherein p is 1 or 2 and each R^1 is chloro.

147. (new) A compound of claim 115 wherein p is 1 or 2 and each R^1 is chloro.

148. (new) A compound of claim 116 wherein p is 1 or 2 and each R^1 is chloro.

149. (new) A compound of claim 117 wherein p is 1 or 2 and each R^1 is chloro.

150. (new) A compound of claim 118 wherein p is 1 or 2 and each R^1 is chloro.

151. (new) A compound of claim 119 wherein p is 1 or 2 and each R^1 is chloro.

152. (new) A compound of claim 120 wherein p is 1 or 2 and each R^1 is chloro.

153. (new) A compound of claim 121 wherein p is 1 or 2 and each R^1 is chloro.

154. (new) A compound of claim 122 wherein p is 1 or 2 and each R^1 is chloro.

155. (new) A compound of claim 123 wherein p is 1 or 2 and each R^1 is chloro.

156. (new) A compound of claim 124 wherein p is 1 or 2 and each R^1 is chloro.

157. (new) A compound of claim 125 wherein p is 1 or 2 and each R^1 is chloro.

158. (new) A compound of claim 126 wherein p is 1 or 2 and each R^1 is chloro.

159. (new) A compound of claim 127 wherein p is 1 or 2 and each R^1 is chloro.

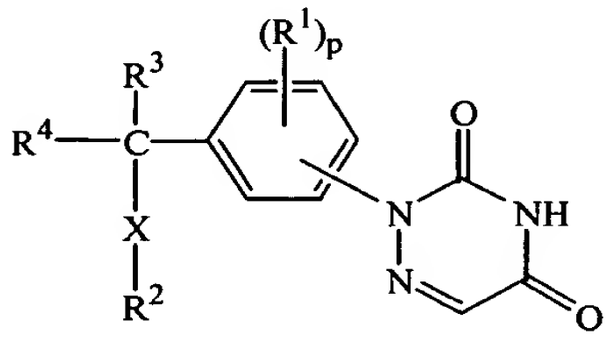
160. (new) A compound of claim 126 wherein p is 1 or 2 and each R^1 is chloro.

161. (new) A compound of claim 111 wherein R^3 and R^4 are both methyl, -X-R^2 is optionally substituted 2-thiazolyl or 3-oxadiazolyl, the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents, and p is 2 whereby

both R^1 substituents are chloro positioned ortho relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

162. (new) A compound of claim 112 wherein R^3 and R^4 are both methyl, -X-R^2 is optionally substituted 2-thiazolyl or 3-oxadiazolyl, the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents, and p is 2 whereby both R^1 substituents are chloro positioned ortho relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

163. (new) A compound having the formula



(I)

a *N*-oxide, a pharmaceutically acceptable addition salt or a stereochemically isomeric form thereof, wherein :

p represents an integer being 0, 1 or 2;

X represents O, S, NR^5 or a direct bond;

Y represents O, S, NR^5, or $S(O)_2$;

each R^1 independently represents chloro or trifluoromethyl;

R^2 represents Het^1 or C_{1-6}alkyl substituted with one or two substituents selected from hydroxy, cyano, amino, mono- or di(C_{1-4}alkyl)amino, C_{1-6}alkyloxy, C_{1-6}alkylsulfonyloxy, C_{1-6}alkyloxycarbonyl, C_{3-7}cycloalkyl, aryl, aryloxy, arylthio, Het^1, Het^1oxy and Het^1thio; and if X is O, S or NR^5, then R^2 may also represent aminocarbonyl, aminothiocarbonyl, C_{1-4}alkylcarbonyl, C_{1-4}alkylthiocarbonyl, arylcarbonyl, arylthiocarbonyl, Het^1carbonyl or Het^1thiocarbonyl;

R^3 represents hydrogen, C_{1-6}alkyl or C_{3-7}cycloalkyl;

R^4 represents hydrogen, C_{1-6}alkyl or C_{3-7}cycloalkyl; or R^3 and R^4 taken together form a C_{2-6}alkanediyl;

R^5 represents hydrogen or C_{1-4}alkyl;

each R^6 independently represents C_{1-6}alkylsulfonyl, aminosulfonyl, mono- or di(C_{1-4}alkyl)aminosulfonyl, mono- or di(benzyl)aminosulfonyl, polyhaloC_{1-6}alkylsulfonyl, C_{1-6}alkylsulfinyl, phenylC_{1-4}alkylsulfonyl, piperazinylsulfonyl, aminopiperidinylsulfonyl, piperidinylaminosulfonyl, *N*-C_{1-4}alkyl-*N*-piperidinylaminosulfonyl or mono-or di(C_{1-4}alkyl)aminoC_{1-4}alkylsulfonyl;

each R^7 and each R^8 are independently selected from hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, aryl, arylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, aminocarbonyl, arylcarbonyl, Het^3carbonyl, C_{1-4}alkylcarbonyloxy-C_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)-aminoC_{1-4}alkyl, arylaminocarbonyl, arylaminothiocarbonyl, Het^3aminocarbonyl, Het^3aminothiocarbonyl, C_{3-7}cycloalkyl, pyridinylC_{1-4}alkyl, C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3, Het^4 and R^6;

R^9 and R^{10} are each independently selected from hydrogen, C_{1-4}alkyl, hydroxyC_{1-4}alkyl, dihydroxyC_{1-4}alkyl, phenyl, phenylC_{1-4}alkyl, C_{1-4}alkyloxyC_{1-4}alkyl, C_{1-4}alkylcarbonyl, aminocarbonyl, phenylcarbonyl, Het^3carbonyl, C_{1-4}alkylcarbonyloxyC_{1-4}alkylcarbonyl, hydroxyC_{1-4}alkylcarbonyl, C_{1-4}alkyloxycarbonylcarbonyl, mono- or di(C_{1-4}alkyl)-aminoC_{1-4}alkyl, phenylaminocarbonyl, phenylaminothiocarbonyl, Het^3aminocarbonyl, Het^3aminothiocarbonyl, C_{3-7}cycloalkyl, pyridinylC_{1-4}alkyl, C_{1-4}alkanediyl-C(=O)-O-R^{14}, -C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3, Het^4 and R^6;

each R^{11} independently being selected from hydroxy, cyano, nitro, halo, C_{1-4}alkyloxy, formyl, NR^7R^8, $C(=O)NR^7R^8$, -C(=O)-O-R^{14}, aryl, arylcarbonyl, Het^3 and $C(=O)Het^3$;

each R^{14} independently represents hydrogen, C_{1-4}alkyl, C_{3-7}cycloalkyl, aminocarbonylmethylene or mono-or di(C_{1-4}alkyl)aminocarbonylmethylene;

aryl represents phenyl optionally substituted with one, two or three substituents each independently selected from nitro, azido, cyano, halo, hydroxy, C_{1-4}alkyl, C_{3-7}cycloalkyl, C_{1-4}alkyloxy, formyl, polyhaloC_{1-4}alkyl, NR^9R^{10}, $C(=O)NR^9R^{10}$, C(=O)-O-R^{14}, R^6, -O-R^6, phenyl, Het^3, $C(=O)Het^3$ and C_{1-4}alkyl substituted with hydroxy, C_{1-4}alkyloxy, C(=O)-O-R^{14}, -Y-C_{1-4}alkanediyl-C(=O)-O-R^{14}, Het^3 or NR^9R^{10};

Het^1 represents a heterocycle selected from pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}; provided Het^1 is other than 2-substituted-pyridin-5-yl;

Het^2 represents a heterocycle selected from furanyl, thienyl, pyridinyl or benzothienyl, wherein said aromatic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^4, R^{11} and C_{1-4}alkyl optionally substituted with R^{11};

Het^3 represents a monocyclic heterocycle selected from piperidinyl, piperazinyl, morpholinyl and tetrahydropyranyl each independently and optionally substituted with, where possible, one, two, three or four substituents each independently selected from hydroxy, C_{1-4}alkyl, C(=O)-O-R^{14}, C_{1-4}alkylcarbonyl, R^6, piperidinyl and C_{1-4}alkyl substituted with one or two substituents independently selected from hydroxy, C_{1-4}alkyloxy, C(=O)-O-R^{14} and phenyl;

Het^4 represents a monocyclic heterocycle selected from thienyl or pyridinyl.

164. (new) A compound of claim 163, wherein when X is a direct bond, at least one of R^3 and R^4 is hydrogen, and R^2 is 3-pyridinyl, then R^2 is not substituted in the 6 position with an optionally substituted alkyl or acyl group.

165. (new) A compound of claim 163 wherein the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.

166. (new) A compound of claim 163 wherein R^2 is a monocyclic heterocycle selected from the group consisting of: pyrrolyl, imidazolyl, pyrazolyl, triazolyl, tetrazolyl, furanyl, thienyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, thiadiazolyl, oxadiazolyl, pyridinyl, pyrimidinyl, pyrazinyl, pyranyl, pyridazinyl and triazinyl, wherein said monocyclic heterocycles each independently may optionally be substituted with one, or where possible, two or three substituents each independently selected from Het^2, R^{11} and C_{1-4}alkyl optionally substituted with Het^2 or R^{11}.

167. (new) A compound of claim 163 wherein R^3 and R^4 are both methyl and -X-R^2 is Het^1.

168. (new) A compound of claim 163 wherein p is 1 or 2 and each R^1 is chloro.

169. (new) A compound of claim 163 wherein R^3 and R^4 are both methyl, -X-R^2 is optionally substituted 2-thiazolyl or 3-oxadiazolyl, the 6-azauracil moiety is in the para position relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents, and p is 2 whereby both R^1 substituents are chloro positioned ortho relative to the carbon atom bearing the -X-R^2, R^3 and R^4 substituents.